

## globals

Defines input namelists and global variables

## contents

<b>1</b>	<b>globals</b>	<b>1</b>
1.1	overview	1
1.2	input list: <code>focusin</code>	1

### 1.1 overview

- Here, and elsewhere, input variables are shown in **red**. The input list is read from file and broadcast in **al00aa**.

### 1.2 input list: `focusin`

- **`Idisplay = 0`** : silent output; -1 = details ;
- **`Isymmetric = 1`** : enforce stellarator symmetry;
- **`Itopology = 0`** : selects knottedness of plasma:
- **`knotsurf = 0.200D-00`** : radius of knotted plasma boundary;
- **`ellipticity = 0.000D-00`** : radius of knotted plasma boundary;
- **`Linitialize = 0`** :
- **`Rmaj = 1.000D+00`** : major radius of coils;
- **`rmin = 0.500D+00`** : minor radius of coils;
- **`Ic = 0`** :
- **`Io = 1.000D+00`** :
- **`Iw = 1.000D+00`** :
- **`Lc = 0`** : logical flag controlling length weight; see **`tlength`**;
- **`Lo = 1.000D+00`** :
- **`Lw = 1.000D+00`** :
- **`NFcoil = 4`** : Fourier harmonics for each coil;
- **`NDcoil = 128`** : discrete segments per coil;
- **`Loptimize = -2/-1/0/1/2/3`** : -1 and -2 are for testing the derivatives; 1 old descent; 2 new descent; 3 Powell nonlinear equations slover; 4 Newton;
- **`Lnormalize = 0/1`** : turn off/on normalizing weights;
- **`weight.bnrm = 1.000D+00`** : weight for bnormal constraint; **`bnormal`**
- **`weight.tflux = 0.500D+00`** : weight for toroidal flux constraint; **`torflux`**
- **`target.tflux = 1.000D+00`** : target toroidal flux; **`torflux`**
- **`weight.ttlen = 0.000D+00`** : weight for coil length; **`tlength`**
- **`weight.eqarc = 1.000D+00`** : weight for equal arc length constraint; **`equarcl`**
- **`weight.ccsep = 0.000D+00`** : weight for coil-coil separation ; **`coilsep`**
- **`tauend = 1.000D-00`** : artificial relaxtion “time”, **`evolve`**;
- **`tautol = 1.000D-04`** : o.d.e. integration tolerance;
- **`Ntauout = 100`** : intermediate time steps; **`evolve`**;

- **Savfreq = 1** : Saving Frequency; 1 means saving files for each step;
- **Nteta = 64** :
- **Nzeta = 64** :
- **absacc = 1.000D-08** :
- **absreq = 1.000D-12** :
- **relreq = 1.000D-01** :
- **xtol = 0.000D+00** : E04LBF tolerance  $10 \times \text{sqrtmachprec}$ ;
- **eta = 0.900D+00** : E04LBF accuracy rate (step ration);
- **stepmx = 1.000D+05** : E04LBF Euclidean distance between solution and starting;
- **Mpol = -8** : Fourier poloidal resolution;
- **Ntor = 4** : Fourier toroidal resolution;
- **Lpoincare = 0** : to construct Poincaré plot;
  - if **Lpoincare** > 0, then the fieldline parameter is the cylindrical toroidal angle, and so  $B^\phi$  must not equal zero;
  - if **Lpoincare** = -1, then the fieldline parameter is the length;
- **odetol = 1.000D-10** : Poincaré plot, **pp00aa**;
- **Ppts = 100** : Poincaré plot, **pp00aa**;
- **Ptrj = 8** : Poincaré plot, **pp00aa**;
- **phi = 0.0** : REDUNDANT;
- **iphi = 0** : Poincaré plot,
- **bstol = 1.000D-06** : tolerance in Biot-Savart integral; passed to **oculus:bs00aa**;
- **bsnlimit = 100000** : max. number of iterations used in Biot-Savart integral; passed to **oculus:bs00aa**;